

EAST Search History

Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
S1	2	"4067722".pn.	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 10:45
S2	98	terpyridine	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 10:47
S3	1483	terpyridine	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 10:47
S4	192	terpyridine.clm.	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 17:01
S5	1033	tetraphenylporphyrin	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 11:27
S6	0	S5 same paranitro	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 11:25
S7	0	S5 same (p adj no)	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 11:25
S8	23	tetraphenylporphyrin same aminophenyl	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 11:28

EAST Search History

S9	0	"tetra(aminophenyl)porphyrin"	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 11:28
S10	3	tetra adj aminophenyl adj porphyrin	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 11:29
S11	188	crapo.in.	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 12:53
S12	29	S11 and porphyrin	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 14:13
S13	1347	porphyrin same peptide	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 14:13
S14	520	porphyrin same peptide same conjugate	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 14:13
S15	2	wo-9414440-\$.did.	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 14:44
S16	0	wo-2001007442-\$.did.	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 14:44

EAST Search History

S17	1	wo-200107442-\$.did.	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 15:14
S18	2	"20040029181".pn.	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 15:14
S19	1	S18 and cleaved	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 15:14
S20	1	1994-234323.NRAN.	DERWENT	AND	ON	2007/12/20 16:07
S21	10	"1359901"	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 16:09
S22	1	2002-644653.NRAN.	DERWENT	AND	ON	2007/12/20 16:10
S23	8	Ru adj polypyridine adj complex	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 16:28
S24	72	polypyridine adj complex	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 16:28
S25	644	terpyridine same complex	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 17:01
S26	477	terpyridine same complex same metal	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 17:02

EAST Search History

S27	36	terpyridine same complex same metal.clm.	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 17:08
S28	2	"5994339".pn.	US-PGPUB; USPAT; USOCR; FPRS; EPO; JPO; DERWENT	AND	ON	2007/12/20 17:08

=> d his ful

(FILE 'HOME' ENTERED AT 09:34:44 ON 20 DEC 2007)

FILE 'REGISTRY' ENTERED AT 09:34:53 ON 20 DEC 2007

L1 STRUCTURE UPLOADED
 D
L2 2 SEA SSS SAM L1
L3 27 SEA SSS FUL L1
L4 ANALYZE PLU=ON L3 1- LC : 7 TERMS
 D

FILE 'ZCAPLUS' ENTERED AT 09:36:18 ON 20 DEC 2007

L5 29 SEA ABB=ON PLU=ON L3
 D L5 1-29 IBIB HITSTR
 D L4

FILE 'BEILSTEIN' ENTERED AT 09:38:21 ON 20 DEC 2007

L6 1 SEA ABB=ON PLU=ON L3
 D
 D L4

FILE 'CHEMCATS, TOXCENTER' ENTERED AT 09:39:03 ON 20 DEC 2007

L7 4 SEA ABB=ON PLU=ON L3
 D L7 1-4

FILE HOME

FILE REGISTRY

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 19 DEC 2007 HIGHEST RN 958936-22-6

DICTIONARY FILE UPDATES: 19 DEC 2007 HIGHEST RN 958936-22-6

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TSCA INFORMATION NOW CURRENT THROUGH June 29, 2007

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<http://www.cas.org/support/stngen/stndoc/properties.html>

FILE ZCAPLUS

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FILE COVERS 1907 - 20 Dec 2007 VOL 147 ISS 26
FILE LAST UPDATED: 19 Dec 2007 (20071219/ED)

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This file contains CAS Registry Numbers for easy and accurate substance identification.

FILE BEILSTEIN
FILE LAST UPDATED ON September 26, 2007

FILE COVERS 1771 TO 2007.
FILE CONTAINS 10.119,480 SUBSTANCES

>>>PLEASE NOTE: Reaction Data and substance data are stored in separate documents and can not be searched together in one query. Reaction data for BEILSTEIN compounds may be displayed immediately with the display codes PRE (preparations) and REA (reactions). A substance answer set retrieved after the search for a chemical name, a compounds with available reaction information by combining with PRE/FA, REA/FA or more generally with RX/FA. The BEILSTEIN Registry Number (BRN) is the link between a BEILSTEIN compound and belonging reactions. For more detailed reaction searches BRNs can be searched as reaction partner BRNs Reactant BRN (RX.RBRN) or Product BRN (RX.PBRN).<<<

>>> FOR SEARCHING PREPARATIONS SEE HELP PRE <<<

* PLEASE NOTE THAT THERE ARE NO FORMATS FREE OF COST. *
* SET NOTICE FEATURE: THE COST ESTIMATES CALCULATED FOR SET NOTICE *
* ARE BASED ON THE HIGHEST PRICE CATEGORY. THEREFORE; THESE *
* ESTIMATES MAY NOT REFLECT THE ACTUAL COSTS. *
* FOR PRICE INFORMATION SEE HELP COST *

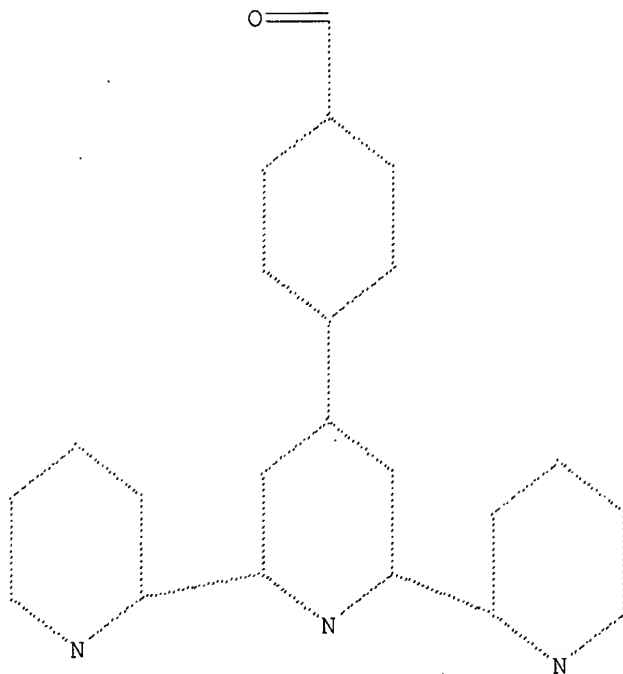
>>> Price change as of January 1st, 2008: Connect Time and Structure Search fees re-introduced. See NEWS and HELP COST <<<

FILE CHEMCATS
FILE LAST UPDATED 15 DECEMBER 2007 (20071215/UP)

FILE TOXCENTER
FILE COVERS 1907 TO 18 Dec 2007 (20071218/ED)

=> d que sta

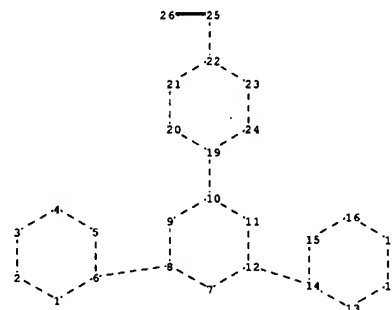
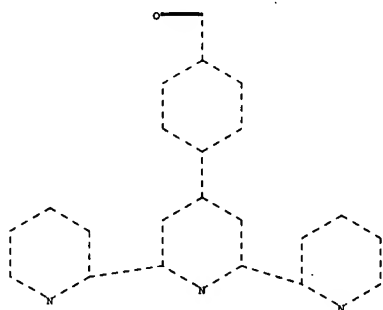
L1 STR



Structure attributes must be viewed using STN Express query preparation.

L3 27 SEA FILE=REGISTRY SSS FUL L1

L7 4 SEA L3



chain nodes :

25 26

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21
22 23 24

chain bonds :

6-8 10-19 12-14 22-25 25-26

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 7-8 7-12 8-9 9-10 10-11 11-12 13-14
13-18 14-15 15-16 16-17 17-18 19-20 19-24 20-21 21-22 22-23 23-24

exact/norm bonds :

1-2 1-6 2-3 3-4 4-5 5-6 6-8 7-8 7-12 8-9 9-10 10-11 10-19
11-12 12-14 13-14 13-18 14-15 15-16 16-17 17-18 19-20 19-24 20-21
21-22 22-23 22-25 23-24 25-26

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom
10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom
18:Atom 19:Atom 20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:CLASS
26:CLASS

=> d his ful

(FILE 'HOME' ENTERED AT 14:22:59 ON 20 DEC 2007)

FILE 'REGISTRY' ENTERED AT 14:25:18 ON 20 DEC 2007

L1 STRUCTURE UPLOADED
L2 STRUCTURE UPLOADED
L3 STRUCTURE UPLOADED
L4 49 SEA SSS SAM L1
L5 50 SEA SSS SAM L2
L6 1008 SEA SSS FUL L1
L7 1336 SEA SSS FUL L2
L8 244 SEA SUB=L6 SSS FUL L3
L9 152 SEA SUB=L7 SSS FUL L3
L10 STRUCTURE UPLOADED
L11 102 SEA SUB=L8 SSS FUL L10
L12 77 SEA SUB=L9 SSS FUL L10
L13 179 SEA ABB=ON PLU=ON L11 OR L12
L14 ANALYZE PLU=ON L13 1- LC : 6 TERMS
 D

FILE 'ZCAPLUS' ENTERED AT 14:30:51 ON 20 DEC 2007

L15 57 SEA ABB=ON PLU=ON L13

FILE 'REGISTRY' ENTERED AT 14:32:12 ON 20 DEC 2007

L16 STRUCTURE UPLOADED
L17 0 SEA SUB=L13 SSS FUL L16

FILE 'ZCAPLUS' ENTERED AT 14:32:57 ON 20 DEC 2007

D L15 1-57 IBIB HITSTR

FILE HOME

FILE REGISTRY

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

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FILE ZCAPLUS

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FILE LAST UPDATED: 19 Dec 2007 (20071219/ED)

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=> d que sta l15

L1 STR

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

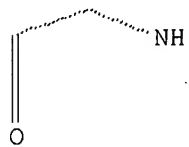
Structure attributes must be viewed using STN Express query preparation.

L2 STR

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

Structure attributes must be viewed using STN Express query preparation.

L3 STR



Structure attributes must be viewed using STN Express query preparation.

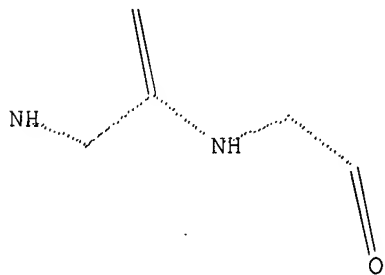
L6 1008 SEA FILE=REGISTRY SSS FUL L1

L7 1336 SEA FILE=REGISTRY SSS FUL L2

L8 244 SEA FILE=REGISTRY SUB=L6 SSS FUL L3

L9 152 SEA FILE=REGISTRY SUB=L7 SSS FUL L3

L10 STR



Structure attributes must be viewed using STN Express query preparation.

L11 102 SEA FILE=REGISTRY SUB=L8 SSS FUL L10

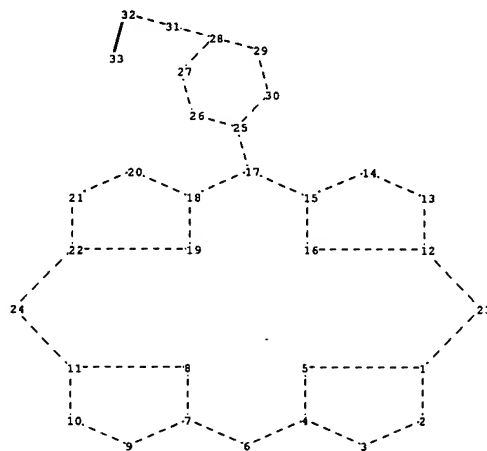
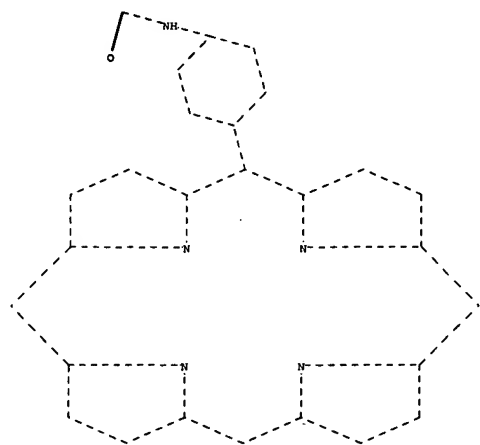
L12 77 SEA FILE=REGISTRY SUB=L9 SSS FUL L10

10/614,324

12/20/2007

L13
L15

179 SEA FILE=REGISTRY ABB=ON PLU=ON L11 OR L12
57 SEA FILE=ZCAPLUS ABB=ON PLU=ON L13



chain nodes :

31 32 33

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21
22 23 24 25 26 27 28 29 30

chain bonds :

17-25 28-31 31-32 32-33

ring bonds :

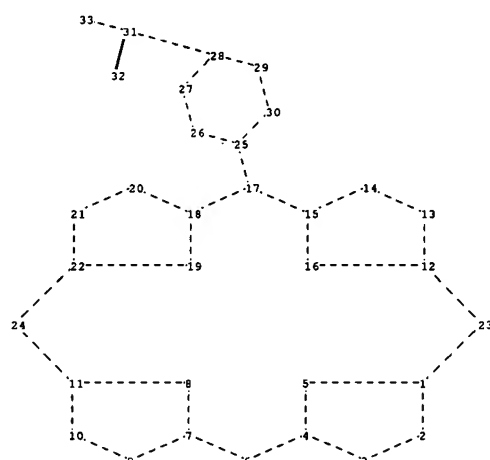
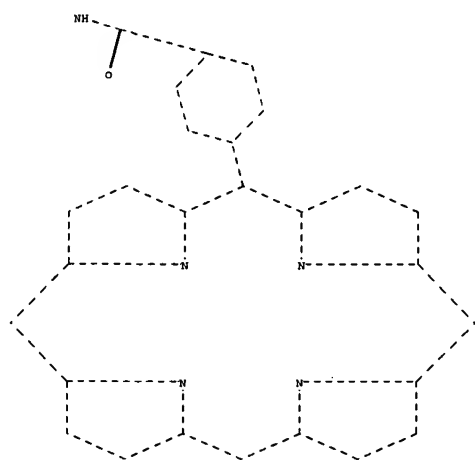
1-2 1-5 1-23 2-3 3-4 4-5 4-6 6-7 7-8 7-9 8-11 9-10 10-11 11-24
12-13 12-16 12-23 13-14 14-15 15-16 15-17 17-18 18-19 18-20 19-22
20-21 21-22 22-24 25-26 25-30 26-27 27-28 28-29 29-30

exact/norm bonds :

1-2 1-5 1-23 2-3 3-4 4-5 4-6 6-7 7-8 7-9 8-11 9-10 10-11 11-24
12-13 12-16 12-23 13-14 14-15 15-16 15-17 17-18 17-25 18-19 18-20
19-22 20-21 21-22 22-24 25-26 25-30 26-27 27-28 28-29 28-31 29-30
31-32 32-33

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom
10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom
18:Atom 19:Atom 20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:Atom
26:Atom 27:Atom 28:Atom 29:Atom 30:Atom 31:CLASS 32:CLASS 33:CLASS



chain nodes :

31 32 33

ring nodes :

 1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21
 22 23 24 25 26 27 28 29 30

chain bonds :

17-25 28-31 31-32 31-33

ring bonds :

 1-2 1-5 1-23 2-3 3-4 4-5 4-6 6-7 7-8 7-9 8-11 9-10 10-11 11-24
 12-13 12-16 12-23 13-14 14-15 15-16 15-17 17-18 18-19 18-20 19-22
 20-21 21-22 22-24 25-26 25-30 26-27 27-28 28-29 29-30

exact/norm bonds :

 1-2 1-5 1-23 2-3 3-4 4-5 4-6 6-7 7-8 7-9 8-11 9-10 10-11 11-24
 12-13 12-16 12-23 13-14 14-15 15-16 15-17 17-18 17-25 18-19 18-20
 19-22 20-21 21-22 22-24 25-26 25-30 26-27 27-28 28-29 28-31 29-30
 31-32 31-33

Match level :

 1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom
 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom
 18:Atom 19:Atom 20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:Atom
 26:Atom 27:Atom 28:Atom 29:Atom 30:Atom 31:CLASS 32:CLASS 33:CLASS

H:\STN queries\amino acid generic.str



chain nodes :

1 2 3 4

chain bonds :

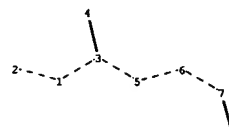
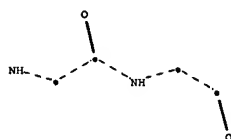
1-3 1-2 3-4

exact/norm bonds :

1-3 1-2 3-4

Match level :

1:CLASS 2:CLASS 3:CLASS 4:CLASS



chain nodes :

1 2 3 4 5 6 7 8

chain bonds :

1-2 1-3 3-4 3-5 5-6 6-7 7-8

exact/norm bonds :

1-2 1-3 3-4 3-5 5-6 6-7 7-8

Match level :

1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS 8:CLASS

H:\STN\queries\generic metal.str

--	--

chain nodes :
1

Match level :
1:CLASS

=> d his ful

(FILE 'HOME' ENTERED AT 17:35:21 ON 20 DEC 2007)

FILE 'REGISTRY' ENTERED AT 17:35:42 ON 20 DEC 2007

L1 STRUCTURE UPLOADED

D

L2 STRUCTURE UPLOADED

L3 1336 SEA SSS FUL L1

L4 152 SEA SUB=L3 SSS FUL L2

FILE 'ZCAPLUS' ENTERED AT 17:37:31 ON 20 DEC 2007

L5 48 SEA ABB=ON PLU=ON L4

D L5 1-48 IBIB HITSTR

FILE 'STNGUIDE' ENTERED AT 17:38:21 ON 20 DEC 2007

FILE HOME

FILE REGISTRY

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FILE STNGUIDE
FILE CONTAINS CURRENT INFORMATION.
LAST RELOADED: Dec 18, 2007 (20071218/UP).

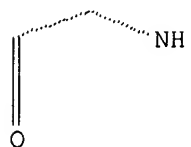
=> d que sta

L1 STR

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

Structure attributes must be viewed using STN Express query preparation.

L2 STR



Structure attributes must be viewed using STN Express query preparation.

L3 1336 SEA FILE=REGISTRY SSS FUL L1
L4 152 SEA FILE=REGISTRY SUB=L3 SSS FUL L2
L5 48 SEA FILE=ZCAPLUS ABB=ON PLU=ON L4